Hypersurfaces and their singularities in partial correlation testing

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Abstract

An asymptotic theory is developed for computing volumes of regions in the parameter space of a directed Gaussian graphical model that are obtained by bounding partial correlations. We study these volumes using the method of real log canonical thresholds from algebraic geometry. Our analysis involves the computation of the singular loci of correlation hypersurfaces. Statistical applications include the strong-faithfulness assumption for the PC-algorithm, and the quantification of confounder bias in causal inference. A detailed analysis is presented for trees, bow-ties, tripartite graphs, and complete graphs.

1 Introduction

Extensive theory has been established in recent years for causal inference based on directed acyclic graph (DAG) models. A popular way for estimating a DAG model from observational data employs partial correlation testing to infer the conditional independence relations in the model. In this paper, we apply algebraic geometry and singularity theory to analyze partial correlations in the Gaussian case. The objects of our study are algebraic hypersurfaces in the parameter space of a given graph that encode conditional independence statements.

We begin with definitions for graphical models in statistics. A DAG is a pair G = (V, E) consisting of a set V of nodes and a set E of directed edges with no directed cycle. We usually take $V = \{1, 2, ..., p\}$ and we associate random variables $X_1, X_2, ..., X_p$ with the nodes. Directed edges are denoted by (i, j) or $i \to j$. The skeleton of a DAG G is the underlying undirected graph obtained by removing the arrowheads. A node i is an ancestor of j if there is a directed path $i \to \cdots \to j$, and a configuration $i \to k, j \to k$ is a collider at k. Finally, we assume that the vertices are topologically ordered, that is $(i, j) \in E$ implies i < j.

Every DAG G specifies a Gaussian graphical model as follows. The adjacency matrix A_G is the strictly upper triangular matrix whose entry in row i and column j is a parameter a_{ij} if $(i,j) \in E$ and it is zero if $(i,j) \notin E$. The Gaussian graphical model is defined by the structural equation model $X = A_G^T X + \epsilon$, where $X = (X_1, \ldots, X_p)^T$. We assume that $\epsilon \sim \mathcal{N}(0, I)$, where I is the $p \times p$ -identity matrix. Then the concentration matrix of this model equals

$$K = (A_G - I)(A_G - I)^T.$$

Key words and phrases: causal inference, PC-algorithm, (strong) faithfulness, real log canonical threshold, resolution of singularities, partial correlation, real radical ideal, asymptotics of integrals, almost-principal minor, directed acyclic graph, Gaussian graphical model, algebraic statistics, singular learning theory.

Since $\det(K) = 1$, the covariance matrix $\Sigma = K^{-1}$ is equal to the adjoint of K. The entries of K and Σ are polynomials in the parameters a_{ij} . Our parameter space for this DAG model will always be a full-dimensional subset Ω of $\mathbb{R}^{|E|}$.

For any subset $S \subset V$ and distinct elements i, j in $V \setminus S$, we represent the conditional independence statement $i \perp \!\!\!\perp j \mid S$ by an almost-principal minor of either K or Σ . By this we mean a square submatrix whose sets of row and column indices differ in exactly one element. To be precise, $i \perp \!\!\!\perp j \mid S$ holds for the multivariate normal distribution with concentration matrix K if and only if the submatrix $K_{iR,jR}$ is singular, where $R = V \setminus (S \cup \{i,j\})$ and $iR = \{i\} \cup R$. The determinant $\det(K_{iR,jR})$ is a polynomial in $(a_{ij})_{(i,j)\in E}$. We are interested in the hypersurface in $\mathbb{R}^{|E|}$ defined by the vanishing of this polynomial. Indeed, the partial correlation is equal to the algebraic expression

$$\operatorname{corr}(i,j|S) = \frac{\det(K_{iR,jR})}{\sqrt{\det(K_{iR,iR}) \cdot \det(K_{jR,jR})}}.$$
(1)

- (a) every non-collider on P is not in S,
- (b) every collider on P is in S or an ancestor of a node in S.

If G has no path that d-connects i and j given S, then i and j are d-separated given S, and $\det(K_{iR,jR}) \equiv 0$ as a function of a. The weight of a path P is the product of all edge weights a_{rs} along this path. It was shown in [17, Equation (11)] that the numerator $\det(K_{iR,jR})$ in (1) is a linear combination, as in (5), of the weights of all paths that d-connect i to j given S. Our primary objects of study are the following subsets of the parameter space:

Tube_{$$i,j|S$$} $(\lambda) = \{ \omega \in \Omega : |\operatorname{corr}(i,j|S)| \le \lambda \}.$ (2)

Here $\operatorname{corr}(i,j|S)$ is a function of the parameter ω (denoted $(a_{ij})_{(i,j)\in E)}$ above) in the space $\Omega\subset\mathbb{R}^{|E|}$, λ is a parameter in [0,1], and (i,j,S) is a triple where i and j are d-connected given S. These "tubes" can be seen as hypersurfaces which have been fattened up by a factor which depends on λ and the position on the hypersurface (see Figure 3). The volume of $\operatorname{Tube}_{i,j|S}(\lambda)$ with respect to a given measure $\varphi(\omega) d\omega$ on $\Omega\subset\mathbb{R}^{|E|}$ is represented by the integral

$$V_{i,j|S}(\lambda) = \int_{\text{Tube}_{i,j|S}(\lambda)} \varphi(\omega) d\omega.$$
 (3)

In this paper we study the asymptotics of this integral when the parameter λ is close to 0.

Two applications in statistics are our motivation. The first concerns the strong-faithfulness assumption for algorithms that learn Markov equivalence classes of DAG models by inferring conditional independence relations. The PC-algorithm [16] is a prominent instance. Our set-up is exactly as in [17]. The Gaussian distribution with concentration matrix K is λ strong-faithful to a DAG G if, for any $S \subset V$ and $i, j \notin S$, we have $|\operatorname{corr}(i, j|S)| \leq \lambda$ if and

only if i is d-separated from j given S. We write $V_G(\lambda)$ for the volume of the region in Ω representing distributions that are not λ -strong-faithful. Then $V_G(\lambda)$ is an aggregate of the volumes in (3) for all non-d-separated triples (i, j, S). Zhang and Spirtes [19] proved uniform consistency of the PC-algorithm under the strong-faithfulness assumption with $\lambda \approx 1/\sqrt{n}$, provided the number of nodes p is fixed and sample size $n \to \infty$. In a high-dimensional, sparse setting, Kalisch and Bühlmann [11] require strong-faithfulness with $\lambda \approx \sqrt{\deg(G) \log(p)/n}$, where $\deg(G)$ denotes the maximal degree (i.e., sum of indegree and outdegree) of nodes in G.

For the scenarios described above it is essential to determine the asymptotic behavior of the unfaithfulness volume $V_G(\lambda)$ when λ tends to 0. We shall address this issue using real log canonical thresholds [1, 14, 18]. Our Section 3 establishes the existence of positive constants ℓ, m, C (which depend on G and φ) such that, asymptotically for $\lambda \to 0$,

$$\begin{array}{lcl}
V_G(\lambda) & \approx & C \cdot \lambda^{\ell} \cdot (-\ln \lambda)^{m-1}, \\
V_{i,j|S}(\lambda) & \approx & C' \cdot \lambda^{\ell'} \cdot (-\ln \lambda)^{m'-1}.
\end{array} \tag{4}$$

(See (9) for an exact definition of \approx .) This refines the results in [17] on the growth of $V_G(\lambda)$ via the geometry of the correlation hypersurfaces $\{\det(K_{iR,jR}) = 0\}$. While [17] focused on developing bounds on $V_G(\lambda)$ for the low-dimensional as well as the high-dimensional case and showed the importance of the number and degrees of these hypersurfaces, we here analyze the exact asymptotic behavior of $V_G(\lambda)$ for $\lambda \to 0$ and G fixed and demonstrate the importance of the singularities of these hypersurfaces. Singularities get fattened up much more than smooth parts of the hypersurface, and this increases the volumes (4) substantially.

Our second application concerns stratification bias in causal inference (see e.g. [7, 8]). Here, the volume $V_{i,j|S}(\lambda)$ being large is not a problematic feature, but is in fact desired. Suppose we want to study the effect of an exposure E on a disease outcome D. If there is an additional variable C such that $D \to C \leftarrow E$, then stratifying (i.e. conditioning) on C tends to change the effect of E on D. This can lead to biases in effect estimation. This is known as colliderbias. On the other hand, if $D \leftarrow C \to E$ holds, then C is a confounder and stratifying on C corresponds to bias removal. In certain larger graphs, such as Greenland's bow-tie example [7], stratifying on C removes confounder-bias but at the same time introduces collider-bias. In order to decide whether one should stratify on such a variable C, it is important to understand the partial correlations involved. In this application, the volume $V_{i,j|S}(\lambda)$ can be viewed as the cumulative distribution function of the partial correlation $\operatorname{corr}(i,j|S)$, and one is interested in comparing the two cumulative distribution functions $V_{E,D|C}(\lambda)$ and $V_{E,D}(\lambda)$.

In this paper we examine $V_{i,j|S}(\lambda)$ from a geometric perspective, and we demonstrate how this volume can be calculated using tools from singular learning theory. To derive the asymptotics (4), the main player is the correlation hypersurface, which is the locus in Ω where $\operatorname{corr}(i,j|S)$ vanishes. The first question is whether this hypersurface is smooth, and, if not, one needs to analyze the nature of its singularities. We study these questions for various classes of interesting causal models, using methods from computational algebraic geometry.

The remainder of this paper is organized as follows: In Section 2 we introduce the families of DAGs which we will be working with throughout. Example 2.1 illustrates the algebraic computations that are involved in our analysis. We also discuss some simulation results, which indicate the importance of singularities when studying the volume $V_G(\lambda)$ of strong-unfaithful

distributions. Section 3 presents the connection to singular learning theory [14, 18] and explains how this theory can be used to compute the volumes of the tubes $\text{Tube}_{i,j|S}(\lambda)$. Example 3.1 illustrates our theoretical results for some very simple polynomials in two variables.

In Section 4 we develop algebraic algorithms for analyzing the singularities of the correlation hypersurfaces. We show that, for the polynomials $\det(K_{iR,jR})$ of interest, the real singular locus is often much simpler than the complex singular locus. For instance, Theorem 4.1 states that these hypersurfaces are always smooth for complete DAGs with up to six nodes. In Section 5 we study the singularities and the volumes (3) for trees without colliders.

Section 6 focuses on our second application, namely bias reduction in causal inference. Problems 6.2 and 6.7 offer precise versions of conjectures by Greenland [7], in terms of comparing different volumes $V_{i,j|S}(\lambda)$ for fixed G. We establish some instances of these conjectures.

In Section 7 we introduce more advanced methods, based on the resolution of singularities [9, 10], for finding the exponents ℓ and m in (4). Finally, in Section 8 we present some new results on computing the constants C and C' in our asymptotics (4) for tube volumes.

2 Four classes of graphs

In this article we will be primarily working with four classes of DAGs:

- i) Complete graphs: We denote the complete DAG on p nodes by K_p . The corresponding matrix A_{K_p} is strictly upper triangular and all $\binom{p}{2}$ parameters a_{ij} are present.
- ii) Trees: We call a DAG G a tree graph if the skeleton of G is a rooted tree and all edges point away from the root (i.e. G has no colliders). We are particularly interested in the most extreme trees, namely star and chain graphs. We denote the star graph shown in Figure 1(b) by Star_p and the chain graph shown in Figure 1(b) by Chain_p .
- iii) Complete tripartite graphs: Let $A, B \subset V$ with $A \cap B = \emptyset$. Then we denote by $A \Rightarrow B$ the complete bipartite graph where $(a,b) \in E$ for all $a \in A$ and $b \in B$. A complete tripartite graph is denoted by Tripart_{p,p'} with $1 \leq p' \leq p-3$. It corresponds to the DAG $\{1,2\} \Rightarrow \{3,\ldots,p-p'\} \Rightarrow \{p-p'+1,\ldots,p\}$ and is shown in Figure 1(c).

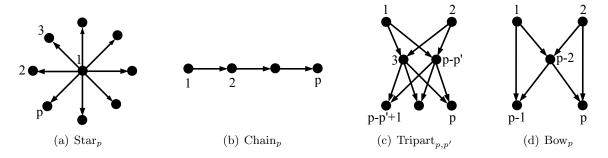


Figure 1: Various classes of graphs.

iv) Bow-ties: We define a bow-tie to be a complete tripartite graph Tripart_{p,2} with two additional edges, namely (1, p - 1) and (2, p). A bow-tie is denoted by Bow_p and is shown in Figure 1(d). Bow-ties with p = 5 feature prominently in Greenland's study [7].

The following example serves as a preview to the topics covered in this paper.

Example 2.1. We illustrate our objects of study for the tripartite graph $G = \text{Tripart}_{6,2}$. This DAG model has eight free parameters, namely the unknowns in the matrix

The covariance matrix Σ equals the inverse (or the adjoint) of the concentration matrix

$$K = \begin{pmatrix} a_{13}^2 + a_{14}^2 + 1 & a_{13}a_{23} + a_{14}a_{24} & -a_{13} & -a_{14} & 0 & 0 \\ a_{13}a_{23} + a_{14}a_{24} & a_{23}^2 + a_{24}^2 + 1 & -a_{23} & -a_{24} & 0 & 0 \\ -a_{13} & -a_{23} & a_{35}^2 + a_{36}^2 + 1 & a_{35}a_{45} + a_{36}a_{46} & -a_{35} & -a_{36} \\ -a_{14} & -a_{24} & a_{35}a_{45} + a_{36}a_{46} & a_{45}^2 + a_{46}^2 + 1 & -a_{45} & -a_{46} \\ 0 & 0 & -a_{35} & -a_{45} & 1 & 0 \\ 0 & 0 & -a_{36} & -a_{46} & 0 & 1 \end{pmatrix}.$$

One conditional independence statement of interest is $1 \perp \!\!\! \perp 2 \mid \{5,6\}$. Its correlation hypersurface in \mathbb{R}^8 is defined by the almost-principal minor in Σ with rows 156 and columns 256, or the almost-principal minor in K with rows 134 and columns 234. That determinant equals

$$f = (1+a_{46}^2)a_{13}a_{23}a_{35}^2 + (1+a_{45}^2)a_{13}a_{23}a_{36}^2 + (1+a_{35}^2)a_{14}a_{24}a_{46}^2 + (1+a_{36}^2)a_{14}a_{24}a_{45}^2 + a_{13}a_{24}a_{35}a_{45} + a_{13}a_{24}a_{36}a_{46} + a_{14}a_{23}a_{35}a_{45} + a_{14}a_{23}a_{36}a_{46} - 2a_{13}a_{23}a_{35}a_{36}a_{45}a_{46} - 2a_{14}a_{24}a_{35}a_{36}a_{45}a_{46}.$$
 (5)

This is a weighted sum of all paths which d-connect nodes 1 and 2 given $\{5,6\}$. The first term in the formula (5) for $f = \det(K_{134,234})$ corresponds to the path $1 \to 3 \to 5 \leftarrow 3 \leftarrow 2$ in $G = \operatorname{Tripart}_{6,2}$, and the last term corresponds to the path $1 \to 4 \to 5 \leftarrow 3 \to 6 \leftarrow 4 \leftarrow 2$.

Let φ be the Lebesgue probability measure on the cube $\Omega = [-1, +1]^8$. The expression $V_{1,2|56}(\lambda)$ defined in (3) is the volume of the region of parameters $a \in \Omega$ that satisfy

$$|\operatorname{corr}(1, 2 \mid 5, 6)| = \left| \frac{f(a)}{\sqrt{\det(K_{134, 134})} \sqrt{\det(K_{234, 234})}} \right| \le \lambda.$$

As a function in λ , the volume $V_{1,2|56}(\lambda)$ is a cumulative distribution function on $[0,\infty)$. Our aim in this article is to determine the asymptotics of such a function for $\lambda \to 0$.

In Section 3 we shall explain the form of the asymptotics that is promised in (4). In order to find the exponents ℓ and m, the first step is to run the algebraic algorithm in Section 4.

This answers the question whether the hypersurface in Ω defined by f = 0 has any singular points. The set of such points, known as the *singular locus*, is the zero set in Ω of the ideal

$$J = \langle f, \frac{\partial f}{\partial a_{13}}, \frac{\partial f}{\partial a_{14}}, \frac{\partial f}{\partial a_{23}}, \frac{\partial f}{\partial a_{24}}, \frac{\partial f}{\partial a_{35}}, \frac{\partial f}{\partial a_{36}}, \frac{\partial f}{\partial a_{45}}, \frac{\partial f}{\partial a_{46}} \rangle.$$

The tools of Section 4 reveal that its real radical [15] is the intersection of three prime ideals:

$$\sqrt[\mathbb{R}]{J} = \langle \text{ entries of } \begin{pmatrix} a_{13} & a_{14} \\ a_{23} & a_{24} \end{pmatrix} \cdot \begin{pmatrix} a_{35} & a_{36} \\ a_{45} & a_{46} \end{pmatrix} \rangle$$

$$= \langle a_{13}, a_{14}, a_{23}, a_{24} \rangle \cap \langle a_{35}, a_{36}, a_{45}, a_{46} \rangle \cap \langle 2 \times 2 \text{-minors of } \begin{pmatrix} a_{13} & a_{23} & a_{45} & a_{46} \\ a_{14} & a_{24} & -a_{35} & -a_{36} \end{pmatrix} \rangle.$$

Thus the hypersurface $\{f=0\}$ is singular. Its singular locus decomposes into three irreducible varieties, namely two linear spaces of dimension 4 and one determinantal variety of dimension 5. In Section 6 we return to this example, with focus on a statistical application of the cumulative distribution function $V_{1,2|56}(\lambda)$, and we will then show that (ℓ, m) equals (1, 1). \square

This paper extends the work of Uhler, Raskutti, Bühlmann and Yu in [17] on the geometry of the strong-faithfulness assumption in the PC-algorithm. Upper and lower bounds on the volume $V_G(\lambda)$ of the unfaithful region Tube_G(λ) for the low- as well as the high-dimensional setting were derived in [17, §5]. These bounds involved only the number |E| of parameters and the degrees of the correlation hypersurfaces $\{\det(K_{iR,jR}) = 0\}$. The new insight in the current paper is that singularities are essential for the asymptotic behavior of $V_G(\lambda)$ for $\lambda \to 0$.

What led us to this insight was taking a closer look at the simulation results for trees. In [17, §6.1.1] trees were still treated as one single class. We subsequently examined the difference

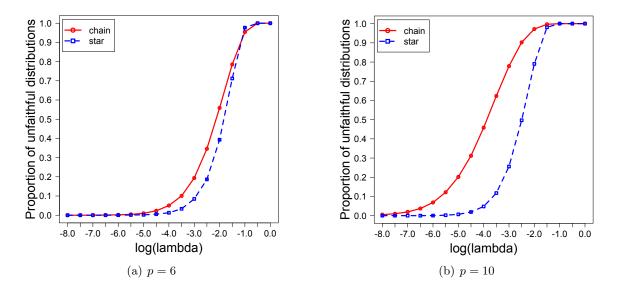


Figure 2: Proportion of λ -strong-unfaithful distributions for chains compared to stars.

between stars and chains, depicted in Figure 1(a) and 1(b). Our simulation results for $Star_p$ and $Chain_p$ are shown in Figure 2. We shall now explain the curves in these diagrams.

The left diagram in Figure 2 is for p = 6 nodes and the right diagram is for p = 10. Each curve is the graph of the cumulative distribution function $V_G(\lambda)$ but with the x-axis transformed to a logarithmic scale (with base 10). Thus we depict the graph of the function

$$(-\infty, 0] \to [0, 1], \ x \mapsto V_G(10^x).$$
 (6)

The red curve is for $G = \operatorname{Chain}_p$ and the blue curve is for $G = \operatorname{Star}_p$. These curves were computed by simulation: we sampled the parameter a from the uniform distribution on $[-1,1]^{p-1}$ and we recorded the proportion of trials that landed in $\operatorname{Tube}_G(\lambda)$ for various values of λ . The diagrams show clearly that $V_G(\lambda)$ is smaller for star graphs than for chain graphs.

A theoretical explanation for these experimental results will be given in Section 5. Our asymptotic theory predicts the behavior of these curves as $x = \log(\lambda)$ tends to $-\infty$. The point is that the correlation hypersurfaces for chain graphs have deeper singularities than those for star graphs. The equation of any such hypersurface for a tree is the product of a monomial and a strictly positive polynomial. This enables us to apply Proposition 3.5. In Theorem 5.1 and Corollary 5.3 we shall determine the constants ℓ , m and m of (4) exactly when the graph m is a tree. We shall also address the question of how to obtain ℓ , m and m from simulations.

Before we get to graphical models, however, we first need to develop the mathematics needed to analyze $V_G(\lambda)$. This will be done, in a self-contained manner, in the next section.

3 Computing the volume of a tube

We now introduce the basics regarding the computation of integrals like the one in (3), and we explain why asymptotic formulas like (4) can be expected. While this section is foundational for what is to follow, no reference to any statistical application is made until Theorem 3.8. It can be read from first principle and might be of independent interest to mathematicians.

Let $\Omega \subset \mathbb{R}^d$ be a compact, full-dimensional, semianalytic subset and consider a probability measure $\varphi(\omega)d\omega$ on Ω where $d\omega$ is the standard Lebesgue measure and $\varphi:\Omega\to\mathbb{R}$ is a real-analytic function. Also, fix an analytic function $f:\Omega\to\mathbb{R}$ whose hypersurface $\{\omega:f(\omega)=0\}$ has non-empty intersection with the interior of Ω . We are interested in the volume $V(\lambda)$ with respect to the measure φ of the region

$$\mathrm{Tube}(\lambda) \ = \ \big\{\omega \in \Omega : |f(\omega)| \le \lambda\big\}.$$

Here $\lambda > 0$ is a parameter that is assumed to be small. In later sections, we often take Ω to be the cube $[-1, +1]^d$, with φ its Lebesgue probability measure, and f is usually a polynomial.

The asymptotics of the volume function $V(\lambda)$ depends on the singularities of the hypersurface $\{f=0\}$. This phenomenon is illustrated in Figure 3. Our measure for the complexity of the singularities of f is a pair (ℓ, m) of non-negative real numbers. That pair is the real log canonical threshold of f. It is related to the volume $V(\lambda)$ for small values of λ by the formula

$$V(\lambda) \approx C \lambda^{\ell} (-\ln \lambda)^{m-1}.$$
 (7)

Here C is a positive real constant whose study we shall defer until Section 8.

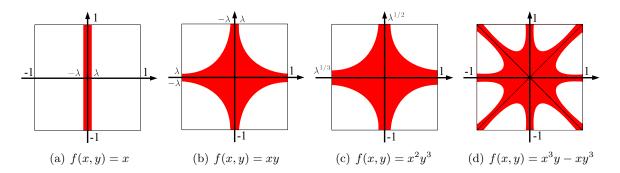


Figure 3: Tubes for various polynomials in two variables.

Example 3.1. Let d=2 and φ the Lebesgue probability measure on the square $\Omega=[-1,+1]^2$. Our problem is to compute the area of the tube $\{(x,y)\in\Omega:|f(x,y)|\leq\lambda\}$. Here f(x,y) is one of the four simple polynomials below whose tubes are shown in Figure 3.

(a) f(x,y) = x: The corresponding tube is a rectangle and its area equals

$$V(\lambda) = \lambda$$
.

So, in this example, we have $(\ell, m) = (1, 1)$ and C = 1. For other lines, the value of C will change. Proposition 3.6 below shows that $(\ell, m) = (1, 1)$ for smooth hypersurfaces.

(b) f(x,y) = xy: The tube in Figure 3(b) consists of four copies of a region that is the union of a small rectangle and a certain area under a hyperbola. Using calculus, we find

$$V(\lambda) = 4\left(\lambda + \int_{\lambda}^{1} \frac{\lambda}{x} dx\right) \frac{1}{4} = \lambda(-\ln \lambda) + \lambda.$$

The logarithm function appears in this case. We have $(\ell, m) = (1, 2)$ and C = 1.

(c) $f(x,y) = x^2y^3$: The corresponding tube is shown in Figure 3(c). Its area equals

$$V(\lambda) \ = \ 4\left(\lambda^{1/2} + \int_{\lambda^{1/2}}^1 \lambda^{1/3} x^{-2/3} dx\right) \frac{1}{4} \ = \ 3\lambda^{1/3} - 2\lambda^{1/2}.$$

So, the real log canonical threshold equals $(\ell, m) = (\frac{1}{3}, 1)$, and we have C = 3. See Proposition 3.5 for a formula for (ℓ, m) when f is a monomial in any number of variables.

(d) f(x,y) = xy(x+y)(x-y): The corresponding tube is shown in Figure 3(d). This example is a slight generalization of (b). As in (b) there is just one singularity at the origin and it is given by the intersection of lines. However, there is no simple closed formula for the area $V(\lambda)$, as this amounts to evaluating a non-trivial abelian integral. We shall see in Example 7.3 that this real log canonical threshold equals $(\ell, m) = (\frac{1}{2}, 1)$.

For general bivariate polynomials f(x,y) we are facing a hard calculus problem, namely integrating the function y=y(x) that is defined implicitly by $f(x,y)=\lambda$. We can approach this by expanding y as a Puiseux series in λ whose coefficients depend on x. Integrating these coefficients leads to asymptotic formulas in λ . These are consistent with what is to follow. \square

We now return to the general setting defined at the beginning of this section. Let W be a random variable taking values in Ω with distribution φ . The volume $V(\lambda)$ with respect to the measure φ can then be viewed as the cumulative distribution function of the random variable f(W). The corresponding probability distribution function $v(\lambda) = dV/d\lambda$ is called the *state density function*. Its Mellin transform is known as the *zeta function* of f. It is denoted by

$$\zeta(z) = \int_0^\infty \lambda^{-z} v(\lambda) d\lambda = \int_\Omega |f(\omega)|^{-z} \varphi(\omega) d\omega \quad \text{for } z \in \mathbb{C}.$$

According to asymptotic theory [1, 14, 18], our volume has the asymptotic series expansion

$$V(\lambda) \approx \sum_{\ell} \sum_{m=1}^{d} C_{\ell,m} \lambda^{\ell} (-\ln \lambda)^{m-1}.$$
 (8)

Here the index ℓ runs over some arithmetic progression of positive rational numbers and d is the dimension of the parameter space Ω . The equation (8) is valid for sufficiently small $\lambda > 0$. To be precise, writing $V(\lambda) \approx \sum_{i=1}^{\infty} g_i(\lambda)$, where $g_1(\lambda) > g_2(\lambda) > \cdots$ for small λ , means that

$$\lim_{\lambda \to 0} \frac{V(\lambda) - \sum_{i=1}^{k} g_i(\lambda)}{g_k(\lambda)} = 0 \quad \text{for each positive integer } k.$$
 (9)

Using the little-o notation, this is equivalent to $V(\lambda) = \sum_{i=1}^k g_i(\lambda) + o(g_k(\lambda))$ as $\lambda \to 0$ for each positive integer k. It is a common misconception to think that the infinite series converges to $V(\lambda)$ for each fixed λ when λ is small. Rather, it means that for each fixed k, the k-term approximation for $V(\lambda)$ gets better as $\lambda \to 0$. We will primarily be interested in the first term approximation (7).

Definition 3.2 ([14, §4.1],[18, §7.1]). We here define the real log canonical threshold (ℓ, m) of f over Ω with respect to φ . This is a pair in $\mathbb{Q}_+ \times \mathbb{Z}_+$ which we denote by $\mathrm{RLCT}_{\Omega}(f; \varphi)$. It measures the complexity of the singularities of the hypersurface defined by $f(\omega) = 0$.

The following four definitions of $RLCT_{\Omega}(f;\varphi) = (\ell,m)$ are known to be equivalent:

(i) For large N > 0, the Laplace integral

$$Z(N) = \int_{\Omega} e^{-N|f(\omega)|} \varphi(\omega) d\omega$$

is asymptotically $CN^{-\ell}(\ln N)^{m-1}$ for some constant C.

(ii) The zeta function

$$\zeta(z) = \int_{\Omega} |f(\omega)|^{-z} \varphi(\omega) d\omega$$

has its smallest pole at $z = \ell$ and that pole has multiplicity m.

(iii) For small $\lambda > 0$, the volume function

$$V(\lambda) = \int_{|f(\omega)| \le \lambda} \varphi(\omega) \, d\omega$$

is asymptotically $C \lambda^{\ell} (-\ln \lambda)^{m-1}$ for some constant C.

(iv) For small $\lambda > 0$, the state density function

$$v(\lambda) = \frac{d}{d\lambda} \int_{|f(\omega)| \le \lambda} \varphi(\omega) \, d\omega$$

is asymptotically $C\lambda^{\ell-1}(-\ln \lambda)^{m-1}$ for some constant C.

If the real analytic hypersurface $\{\omega \in \Omega : f(\omega) = 0\}$ is empty, we set $\ell = \infty$ and we leave m undefined. We order the pairs (ℓ, m) reversely by the size of $\lambda^{\ell}(-\ln \lambda)^{m-1}$ for sufficiently small $\lambda > 0$. This means that $(\ell_1, m_1) < (\ell_2, m_2)$ if $\ell_1 < \ell_2$ or if $\ell_1 = \ell_2$ and $m_1 > m_2$.

Let us provide some intuition for the ordering of the pairs (ℓ, m) . The real log canonical threshold is a measure of complexity for singularities. Analytic varieties can be stratified into subsets where this measure is constant. The highest stratum contains the smooth points of the variety. As we go deeper, to strata with lower real log canonical thresholds, we encounter singularities of increasing complexity. The volumes of λ -fattenings of deeper singularities will also be larger than that of their less complex counterparts. For instance, in Figures 3(b) and 3(c) the singular locus of both examples consists of just the origin, but we observe that the λ -fattening of the origin in Figure 3(c) is larger than in Figure 3(b). See also Example 3.7.

Example 3.3. Let $f(\omega) = \omega_1^2 + \omega_2^2 + \cdots + \omega_d^2$ and φ the Lebesgue probability measure on $\Omega = [-1, -1]^d$. Then Tube(λ) is the *standard ball* of radius $\lambda^{1/2}$, whose φ -volume is

$$V(\lambda) = \frac{\pi^{d/2}}{2^d \cdot \Gamma(\frac{d}{2} + 1)} \cdot \lambda^{d/2}.$$

By Definition 3.2 (c), here the real log canonical threshold equals $RLCT_{\Omega}(f;\varphi) = (d/2,1)$. \square

We now list some formulas for computing the real log canonical threshold. A first useful fact is that $RLCT_{\Omega}(f;\varphi)$ is independent of the underlying measure φ as long as it is positive everywhere. We can thus assume that φ is the uniform distribution on Ω .

Proposition 3.4. If $\varphi : \Omega \to \mathbb{R}$ is strictly positive and 1 denotes the constant unit function on Ω , then

$$RLCT_{\Omega}(f;\varphi) = RLCT_{\Omega}(f;1).$$

Proof. See [14, Lemma 3.8].

Proposition 3.5. Suppose that Ω is a neighborhood of the origin. If $f(\omega) = \omega_1^{\kappa_1} \cdots \omega_d^{\kappa_d} g(\omega)$ where $g: \Omega \to \mathbb{R}$ does not have any real zeros, then $\mathrm{RLCT}_{\Omega}(f; 1) = (\ell, m)$ where

$$\ell = \min_{i} \frac{1}{\kappa_i}$$
 and $m = \left| \underset{i}{\operatorname{argmin}} \frac{1}{\kappa_i} \right|$.

Proof. This is a special case of Theorem 7.1 which will be proved later.

Recall that an analytic hypersurface $\{f(\omega)=0\}$ is singular at a point $\omega \in \Omega$ if ω satisfies

$$f(\omega) = 0$$
 and $\frac{\partial f}{\partial \omega_i}(\omega) = 0$ for $i = 1, \dots, d$.

If the hypersurface is not singular at any point $\omega \in \Omega$, then it is said to be *smooth*.

Proposition 3.6. If the hypersurface $\{f(\omega) = 0\}$ is smooth then $RLCT_{\Omega}(f;1) = (1,1)$.

Proof. This is also a special case of Theorem 7.1.

Example 3.7. Following up on Example 3.1, we now consider an arbitrary monomial function $f(x,y) = x^s y^t$ on the square $\Omega = [-1,1]^2$. The tube looks as in Figure 3(c). Its area satisfies

$$V(\lambda) \approx \begin{cases} C\lambda^{1/s} & \text{if } s < t \\ C\lambda^{1/t} & \text{if } s > t \\ C\lambda^{1/s}(-\ln \lambda) & \text{if } s = t \end{cases}$$

This formula for the asymptotics (7) follows from Definition 3.2 and Proposition 3.5.

For the statistical applications in this paper, the relevant functions f are polynomials. They are determinants $f = \det(K_{iR,jR})$, where $R = V \setminus (S \cup \{i,j\})$ as in Section 1. Let $\mathrm{RLCT}(i,j|S)$ denote the corresponding real log canonical threshold over $\Omega = [-1,1]^E$ with respect to a positive density φ . The theory developed so far says that the real log canonical threshold of the correlation hypersurface gives an asymptotic volume formula for $V_{i,j|S}(\lambda)$.

Theorem 3.8. If φ satisfies the assumptions in Proposition 3.4, then as λ tends to zero, the volume of the region Tube_{i,j|S}(λ) (see (2)) is asymptotically

$$V_{i,j|S}(\lambda) \approx C \lambda^{\ell} (-\ln \lambda)^{m-1}$$

for some constant C > 0 (which only depends on G) and $(\ell, m) = \text{RLCT}(i, j|S)$.

Proof. By part (iii) in Definition 3.2, the desired pair (ℓ, m) is the real log canonical threshold of the partial correlation $f = \operatorname{corr}(i, j|S)$. This is the algebraic (and hence analytic) function in (1). This function differs from the polynomial $\det(K_{iR,jR})$ by a denominator that does not vanish over Ω . That denominator is a unit in the ring of real analytic functions over Ω , and multiplying by a unit does not change the RLCT of an analytic function [14, §4.1].

We close this section by relating our results directly to the study of unfaithfulness in [17].

Corollary 3.9. Under the assumptions in Theorem 3.8, as λ tends to zero, the volume of λ -strong-unfaithful distributions satisfies

$$V_G(\lambda) \approx C \lambda^{\ell} (-\ln \lambda)^{m-1}$$

for some constant C > 0. Here (ℓ, m) is the minimum of the pairs RLCT(i, j|S), where (i, j, S) runs over all triples in the DAG G such that i is not d-separated from j given S.

Proof. The function $V_G(\lambda)$ is the volume of the union of the regions Tube_{i,j|S}(λ). Thus,

$$\max_{i,j,S} V_{i,j|S}(\lambda) \leq V_G(\lambda) \leq \sum_{i,j,S} V_{i,j|S}(\lambda)$$

Asymptotically, for small positive values of λ , both the lower and upper bounds vary like a constant multiple of $\lambda^{\ell}(-\ln \lambda)^{m-1}$ where (ℓ, m) is the minimum over all pairs RLCT(i, j|S). In this minimum, (i, j, S) runs over all triples such that i and j are d-connected given S. \square

4 Singular Locus

The asymptotic integration theory in Section 3 requires us to analyze the singular locus Sing(f) of the real algebraic hypersurface determined by a given polynomial f. If Sing(f) is empty then the hypersurface is smooth and Proposition 3.6 characterizes the asymptotics of the integral. In this section we return to Gaussian graphical models, we develop tools for computing the relevant singular loci, and we show that they are empty in many cases. In many of the remaining cases, the singularities are of the monomial type featured in Proposition 3.5.

Consider any almost-maximal minor $f = \det(K_{iR,jR})$ of the concentration matrix K of a DAG G. This is a polynomial function on the parameter space \mathbb{R}^E . This polynomial and its partial derivatives are elements in the polynomial ring $\mathbb{Q}[a_{ij}:(i,j)\in E]$. The *Jacobian ideal* of f is the ideal in this polynomial ring generated by f and its partials. We denote it by

$$\operatorname{Jacob}_{i,j,R} := \langle f \rangle + \langle \frac{\partial f}{\partial a_{ij}} : (i,j) \in E \rangle.$$

The singular locus $\operatorname{Sing}(f)$ is the subvariety of real affine space \mathbb{R}^E defined by the Jacobian ideal $\operatorname{Jacob}_{i,j,R}$. The structure of the real variety $\operatorname{Sing}(f)$ governs the volume $V_{i,j|S}(\lambda)$ of the set $\operatorname{Tube}_{i,j|S}(\lambda)$ of unfaithful parameters. If $\operatorname{Sing}(f) = \emptyset$ then Proposition 3.6 tells us that $V_{i,j|S}(\lambda)$ asymptotically equals $C\lambda$ for some constant C > 0. If the singular locus is not empty then understanding $\operatorname{Sing}(f)$ is essential for computing its real log canonical threshold (ℓ, m) .

We conducted a comprehensive study of all DAGs with few nodes by computing the singular locus for every almost-principal minor in their concentration matrix K. Our first result concerns the special case of complete graphs. Non-complete graphs will be studied later.

Theorem 4.1. Suppose that φ satisfies the assumptions in Proposition 3.4. For any conditional independence statement on the complete directed graph K_p with $p \leq 6$ nodes, we have $\operatorname{Sing}(f) = \emptyset$, and hence $V_{i,j|S}(\lambda) \approx C\lambda$ for all triples (i,j,S).

It is tempting to conjecture that the hypothesis $p \leq 6$ can be removed in this theorem. Presently we do not know how to approach this problem other than by direct calculation.

Applying Corollary 3.9, this means that the volume of λ -strong-unfaithful distributions for the complete graph satisfies $V_{K_p}(\lambda) \approx C \lambda$ for $\lambda \to 0$, which is the best possible behavior regarding strong-faithfulness. This may be counter-intuitive, but is confirmed in simulations. In Figure 4 we plot (via (6)) the proportion of strong-unfaithful distributions $V_G(\lambda)$ for the five graphs in Section 2 for varying values of λ . Especially in the plot for p=10 it becomes apparent that the behavior for $\lambda \to 0$ is very different than, say, for $\lambda = 0.001$. For $\lambda \to 0$ we have $V_{\text{complete}}(\lambda) < V_{\text{chain}}(\lambda)$, although the chain graph is much sparser than the complete graph. Note also that the complete graph K_{10} has $\sum_{k=2}^{10} {10 \choose k} {k \choose 2} = 11520$ relevant triples (i,j,S), whereas for Chain₁₀ there are only $\sum_{k=1}^{9} k2^{k-1} = 4097$ such triples.

In what follows we explain the algebraic computations that led to Theorem 4.1. We used ideal-theoretic methods from [3] in their implementation in the Gröbner-based software packages Macaulay 2 [5] and Singular [4]. An important point to note at the outset is that the ideal Jacob_{i,j,R} is almost never the unit ideal. By Hilbert's Nullstellensatz, this means that the hypersurfaces $f = \det(K_{iR,jR})$ do have plenty of singular points over the field $\mathbb C$ of

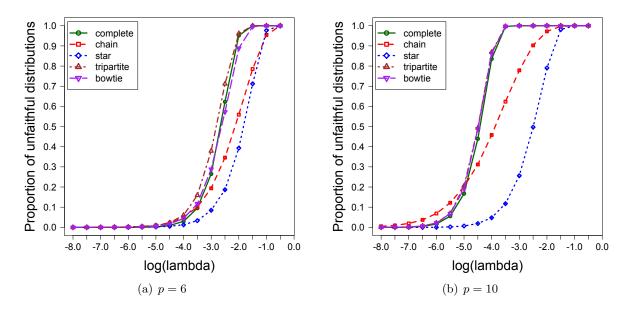


Figure 4: $V_G(\lambda)$ for the complete graph K_p compared to Chain, Star, Tripart, Bow.

complex numbers. What Theorem 4.1 asserts is that, in many of the cases of interest to us here, none of those singular points have their coordinates in the field \mathbb{R} of real numbers.

In order to study the real variety of an ideal, techniques from real algebraic geometry are needed. A key technique is to identify sums of squares (SOS). Indeed, the *real Nullstellensatz* [15] states that the real variety is empty if and only if the given ideal contains a certain type of SOS. To apply this to directed Gaussian graphical models, we shall use the fact that every principal minor of the covariance matrix or the concentration matrix furnishes such an SOS.

Lemma 4.2. Every principal minor $\det(K_{R,R})$ of the concentration matrix K of a DAG is equal to 1 plus a sum of squares in $\mathbb{Q}[a_{ij}:(i,j)\in E]$. In particular, its real variety is empty.

Proof. We can write the principal submatrix $K_{R,R}$ as the product $(A-I)_{R,*} \cdot ((A-I)_{R,*})^T$, where $()_{R,*}$ refers to the submatrix with row indices R. Thus $K_{R,R}$ is the product of an $|R| \times p$ -matrix and its transpose. By the Cauchy-Binet formula, $\det(K_{R,R})$ equals the sum of squares of all maximal minors of the $|R| \times p$ -matrix $(A-I)_{R,*}$. One of these maximal minors is the identity matrix. Hence the polynomial $\det(K_{R,R})$ has the form $1+\mathrm{SOS}$. In particular, the matrix $K_{R,R}$ is invertible for all parameter values in \mathbb{R}^E .

In the context of commutative algebra, it now makes sense to introduce the saturations

$$\begin{array}{rcl} \operatorname{Singu}_{i,j,R} & = & \left(\operatorname{Jacob}_{i,j,R} : \det(K_{R,R})^{\infty}\right), \\ \operatorname{Singu}_{i,j,R}^{*} & = & \left(\operatorname{Singu}_{i,j,R} : \left(\prod_{(i,j)\in E} a_{ij}\right)^{\infty}\right). \end{array}$$

These are also ideals in $\mathbb{Q}[a_{ij}:(i,j)\in E]$. By definition, Singu_{i,j,R} consists of all polynomials that get multiplied into the Jacobian ideal by some power of the determinant of $K_{R,R}$, and Singu^{*}_{i,j,R} consists of polynomials that get multiplied into Singu_{i,j,R} by some monomial. By

[3, §4.4], the variety of $\operatorname{Singu}_{i,j,R}$ is the Zariski closure of the set-theoretic difference of the variety of $\operatorname{Jacob}_{i,j,R}$ and the hypersurface $\{\det(K_{R,R})=0\}$. We saw in Lemma 4.2 that the latter hypersurface has no real points. The ideal $\operatorname{Singu}_{i,j,R}^*$ represents singularities in $(\mathbb{R}^E)^*$.

Corollary 4.3. The singular locus of the real algebraic hypersurface $\{\det(K_{iR,jR}) = 0\}$ in \mathbb{R}^E coincides with the set of real zeros of the ideal $\operatorname{Singu}_{i,j,R}$. The set of real zeros of $\operatorname{Singu}_{i,j,R}^*$ is the Zariski closure of the subset of all singular points whose coordinates are non-zero.

Proof of Theorem 4.1. We computed the ideals $Jacob_{i,j,R}$ and $Singu_{i,j,R}$ for every almost-principal minor $K_{iR,jR}$ in the concentration matrices of the graphs $G = K_3, K_4, K_5, K_6$. In all cases the ideal $Singu_{i,j,R}$ was found to equal the unit ideal $\langle 1 \rangle$. These exhaustive computations were carried out using the software Singular [4]. This establishes Theorem 4.1.

We briefly discuss our computations for the complete directed graph on six nodes.

Example 4.4. Fix the complete directed graph $G = K_6$. We tested all 240 conditional independence statements and computed the corresponding ideal $\operatorname{Singu}_{i,j,R}$. We discuss one interesting instance, namely $i = 1, j = 3, R = \{2, 4\}$. The almost-principal minor $K_{241,243} =$

$$\begin{pmatrix} a_{23}^2 + a_{24}^2 + a_{25}^2 + a_{26}^2 + 1 & a_{25}a_{45} + a_{26}a_{46} - a_{24} & a_{24}a_{34} + a_{25}a_{35} + a_{26}a_{36} - a_{23} \\ a_{25}a_{45} + a_{26}a_{46} - a_{24} & a_{45}^2 + a_{46}^2 + 1 & a_{35}a_{45} + a_{36}a_{46} - a_{34} \\ a_{13}a_{23} + a_{14}a_{24} + a_{15}a_{25} + a_{16}a_{26} - a_{12} & a_{15}a_{45} + a_{16}a_{46} - a_{14} & a_{14}a_{34} + a_{15}a_{35} + a_{16}a_{36} \end{pmatrix}$$

contains all 15 parameters except a_{56} . Its determinant is a polynomial of degree 6. Of its 14 partial derivatives, 13 have degree 5. The derivative with respect to a_{12} has degree 4. Thus $Jacob_{1,3,\{2,4\}}$ is generated by 15 polynomials of degrees $4,5,\ldots,5,6$. The matrix $K_{24,24}$ is the upper left 2×2 -block in the matrix above. The square of its determinant is a polynomial of degree 8 that happens to lie in the ideal $Jacob_{1,3,\{2,4\}}$. This proves $Singu_{1,3,\{2,4\}} = \langle 1 \rangle$.

For graphs G that are not complete, $\operatorname{Singu}_{i,j,R}$ may not be the unit ideal. We already saw one non-obvious instance of this for the tripartite graph in Example 2.1. Here is an even smaller example where the Jacobian ideal and its saturations are equal, and not the unit ideal.

Example 4.5. Let p=4 and take G to be the almost-complete graph with adjacency matrix

$$A_G = \begin{pmatrix} 0 & 0 & a_{13} & a_{14} \\ 0 & 0 & a_{23} & a_{24} \\ 0 & 0 & 0 & a_{34} \\ 0 & 0 & 0 & 0 \end{pmatrix}.$$

The conditional independence statement $1 \perp 1 \leq 1/4$ is represented by the almost-principal minor

$$K_{31,32} = \begin{pmatrix} a_{34}^2 + 1 & a_{24}a_{34} - a_{23} \\ a_{14}a_{34} - a_{13} & a_{13}a_{23} + a_{14}a_{24} \end{pmatrix}$$

of the concentration matrix. The determinant of this minor factors into two binomial factors:

$$\det(K_{31,32}) = (a_{13}a_{34} + a_{14})(a_{23}a_{34} + a_{24}). \tag{10}$$

The Jacobian ideal is the prime ideal generated by these factors:

$$\operatorname{Jacob}_{1,2,3} = \operatorname{Singu}_{1,2,3}^* = \operatorname{Singu}_{1,2,3}^* = \langle a_{13}a_{34} + a_{14}, a_{23}a_{34} + a_{24} \rangle.$$

The left equality holds because $\det(K_{3,3}) = a_{34}^2 + 1$ is a non-zerodivisor modulo $\operatorname{Jacob}_{1,2,3}$. The singular locus of (10) is the three-dimensional real variety defined by this binomial ideal in the parameter space \mathbb{R}^5 . Its real log canonical threshold is found to be $(\ell, m) = (1, 2)$. \square

This example inspired us to analyze the partial correlations of all small DAGs with $p \leq 4$ nodes. In our experiments, we found that $\det(K_{iR,jR})$ is frequently the product of a monomial with a strictly positive sum of squares. Such cases are denoted as "Monomial" in Tables 1 and 2. For these, the RLCT is read off directly from Proposition 3.5. The rows labeled "Smooth" cover cases that are not monomial but where $\operatorname{Singu}_{i,j,R}$ is the unit ideal, so Proposition 3.6 gives us the RLCT. The next theorem summarizes the complete results. The trivial case p=2 is excluded because there is only one graph $1 \to 2$, with $\operatorname{RLCT}(1, 2|\emptyset) = (1, 1)$.

Theorem 4.6. Under the assumptions in Theorem 3.8, for all DAGs with $p \le 4$ nodes and all triples (i, j, S), the value RLCT(i, j|S) is given in Tables 1 and 2. In all cases but one, we have RLCT(i, j|S) = (1, m) where m < p.

To establish Theorem 4.6 we listed every DAG G and every triple (i, j, S) that is not d-separated in G. The rows "Monomial" and "Smooth" were discussed above. The row "Normal crossing" refers to cases that are covered by Theorem 7.1. The "Special" cases are treated in Examples 4.5 and 4.8. Lastly, the row "Blowup" represents instances where the real singular locus is a linear space. Our computation of $\mathrm{RLCT}(i,j|S) = (\ell,m)$ for such instances uses the method explained in Example 7.4. We now examine the unique exceptional case where $\ell \neq 1$.

Table 1: RLCT for all DAGs with three nodes

	(1,1)	(1, 2)	Subtotal
Monomial	21	3	24
Smooth	3		3
Subtotal	24	3	27

Table 2: RLCT for all DAGs with four nodes

	(1, 1)	(1, 2)	(1,3)	(1/2,1)	Subtotal
Monomial	568	145	14	1	728
Smooth	198				198
Normal crossing		22	2		24
Blowup	12				12
Special	2	1			3
Subtotal	780	168	16	1	965

Example 4.7. Let p = 4 and $G = \text{Tripart}_{4,1}$. Its concentration matrix may be obtained from Example 4.5 by setting $a_{14} = a_{24} = 0$. The partial correlation for $1 \perp \!\!\! \perp 2 \mid 4$ is now given by

$$\det(K_{13,23}) = a_{13}a_{23}a_{34}^2.$$

For this monomial, Proposition 3.5 tells us that $(\ell, m) = \text{RLCT}(1, 2|4) = (1/2, 1)$.

Here is an interesting case where the RLCT depends in a subtle way on the choice of Ω .

Example 4.8. Consider the conditional independence statement $1 \perp 1 \mid 3 \mid 4$ for the DAG in Figure 5. The partial correlation is represented by the almost-principal minor

$$\det(K_{12,23}) = a_{13} \cdot g \quad \text{where} \quad g = a_{23}a_{24}a_{34} + a_{24}^2 + 1.$$

The component $\{g=0\}$ is smooth in \mathbb{R}^4 . However, it is disjoint from the cube $\Omega = [-1,1]^4$. To see this, note that $-1 \leq a_{23}a_{24}a_{34}$ in Ω . With this, g=0 would imply $a_{24}=0$ and hence g=1, a contradiction. Consequently, if Ω is the cube $[-1,1]^4$ then the correlation hypersurface is simply $\{a_{13}=0\}$, and the RLCT equals (1,1) by Proposition 3.5. The other special case with RLCT = (1,1) in Table 2 comes from swapping the labels of nodes 1 and 2.

Now, if we enlarge the parameter space Ω then the situation changes. For instance, suppose $(a_{13}, a_{23}, a_{24}, a_{34}) = (0, -2, 1, 1)$ is in the interior of Ω . This is a singular point of $\det(K_{12,23}) = a_{13} \cdot g$. The RLCT can be computed by applying Theorem 7.1. It is now (1,2) instead of (1,1). This example shows that the asymptotics of $V_{i,j|S}(\lambda)$ depends on Ω . \square

Remark 4.9. We briefly return to the issue of faithfulness in the PC-algorithm. Zhang and Spirtes [20] introduced a variant known as the *conservative PC-algorithm*. It is less informative than the PC-algorithm, but only requires adjacency-faithfulness for correct inference. The adjacency-faithfulness condition is simply strong-faithfulness restricted to the edges of G:

$$|\operatorname{corr}(i,j|S)| > \lambda$$
 for all $(i,j) \in E$ and $S \subset V \setminus \{i,j\}$.

If $\{i, j\}$ is not adjacent to R then the relevant minor equals $\det(K_{iR,jR}) = a_{ij}\det(K_{R,R}) + f(\bar{a})$, where f is a polynomial in $\bar{a} = \{a_{st} \mid (s,t) \neq (i,j)\}$, the correlation hypersurface is smooth, and $(\ell, m) = (1, 1)$. If $\{i, j\}$ is adjacent to R then the behavior can be wild, as in Example 4.8.

5 Asymptotics for trees

In [17] trees were treated as one class. However, as already noted when discussing Figure 2, there is a striking difference between the volume $V_G(\lambda)$ for chain graphs compared to stars. In this section, we give an explanation for this difference based on real log canonical thresholds.

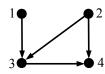


Figure 5: 4-node DAG.

We use the notation SOS(a) for any polynomial that is a sum of squares of polynomials in the model parameters $(a_{ij})_{(i,j)\in E}$. Suppose that G is a tree on $V = \{1, 2, ..., p\}$ and let m be the longest length of an undirected path in G. It was shown in [17, Corollary 4.3(a)] that any non-zero almost-principal minor of the concentration matrix K has the form

$$\det(K_{iR,iR}) = (1 + SOS(a)) \cdot a_{i \to i}, \tag{11}$$

where $a_{i\to j}$ is the monomial of degree $\leq m$ formed by multiplying the parameters a_{rs} along the unique path between i and j. Specifically, for the two trees in Figure 1 we have

$$\det(K_{iR,jR}) = \begin{cases} (1 + SOS(a)) \prod_{k=i}^{j-1} a_{k,k+1} & \text{if } G = \text{Chain}_p, \\ (1 + SOS(a)) \cdot a_{1,i} a_{1,j} & \text{if } G = \text{Star}_p, \text{ and } i, j > 1. \end{cases}$$

In both cases, the term SOS(a) disappears when i and j are leaves of the tree G; cf. (13),(14). Since the correlation hypersurfaces for trees are essentially given by monomials, we can apply Proposition 3.5. The minimal real log canonical threshold is (1, m) where m is the largest degree of any of the monomials in (11). Corollary 3.9 implies the following result:

Theorem 5.1. Under the assumptions in Theorem 3.8, if G is a tree then the volume of λ -strong-unfaithful distributions satisfies

$$V_G(\lambda) \approx C\lambda(-\ln \lambda)^{m-1}$$

where m is the length of the longest path in the tree G, and C is a suitable constant.

For the case of stars we have m=2, whereas for chain graphs we have m=p-1.

Corollary 5.2. Under the assumptions in Theorem 5.1, the volume $V_G(\lambda)$ of strong-unfaithful distributions satisfies

$$V_G(\lambda) \approx \begin{cases} C_{\text{chain}} \cdot \lambda (-\ln \lambda)^{p-2} & \text{if } G = Chain_p, \\ C_{\text{star}} \cdot \lambda (-\ln \lambda) & \text{if } G = Star_p. \end{cases}$$
 (12)

where C_{chain} and C_{star} are suitable positive constants.

As a consequence, the volume $V_G(\lambda)$ is asymptotically larger for chains compared to stars, and the difference increases with increasing number of nodes p. This furnishes an explanation for Figure 2, at least for small values of λ . In that figure we saw the curve for the chain lying clearly above the curve for the star tree. However, one subtle issue is the size of the constants C_{chain} and C_{star} . These need to be understood in order to make accurate comparisons.

In Section 8, we develop new theoretical results regarding the computation of the constant C in (7). Theorem 8.5 gives an integral representation for C when the partial correlation hypersurface is essentially defined by a monomial. In Example 8.7 we shall then derive:

Corollary 5.3. The two constants in (12) are

$$C_{\text{chain}} = \frac{1}{(p-2)!}$$
 and $C_{\text{star}} = \binom{p-1}{2}$.

This result surprised us at first. It establishes the counterintuitive fact that, as p grows, the constant for the lower curve in Figure 2 is exponentially larger than that for the upper curve. Therefore, in order to fully explain the relative position of the two curves for a wider range of values of $\lambda > 0$, it does not suffice to just consider the first order asymptotics (7). Instead, we need to consider some of the higher order terms in the series expansion (8).

As we shall see in Section 8, it is difficult to determine the constants $C_{\ell,m}$ in (8) analytically. In the remainder of this section, we propose a procedure based on simulation and linear regression for estimating the constants $C_{\ell,m}$ in the asymptotic explanations of the volumes $V_G(\lambda)$ or $V_{i,j|S}(\lambda)$. For simplicity we focus on the latter case and we take $f = \det(K_{iR,jR})$.

Suppose that G is a DAG for which the real log canonical thresholds (ℓ, m) in Theorem 3.8 and Corollary 3.9 are known. This is the case for all trees by Theorem 5.1. Our procedure goes as follows. We first sample n points uniformly from Ω and compute the proportion of points ω that lie in Tube_{i,j|S}(λ) for different values of λ . We then fit a linear model to

$$\frac{V_{i,j|S}(\lambda)}{\lambda \ell} \approx C_{m-1}(-\ln \lambda)^{m-1} + C_{m-2}(-\ln \lambda)^{m-2} + \dots + C_0,$$

where (ℓ, m) is the known real log canonical threshold.

In the following, we illustrate this procedure for chain graphs and stars. We analyze two specific examples of partial correlation volumes, namely the ones corresponding to the longest paths in each graph, that is $V_{1,p|\emptyset}(\lambda)$ for Chain_p and $V_{2,3|\emptyset}(\lambda)$ for Star_p. For chain graphs,

$$\operatorname{corr}(1,p) = \frac{\prod_{i=1}^{p-1} a_{i,i+1}}{\sqrt{1 + a_{p-1,p}^2 \left(1 + a_{p-2,p-1}^2 \left(\dots \left(1 + a_{12}^2\right)\right)\right)}},$$
(13)

whereas for star graphs,

$$corr(2,3) = \frac{a_{12} a_{13}}{\sqrt{(1+a_{12}^2)(1+a_{13}^2)}}.$$
(14)

We first approximate $V_{1,p|\emptyset}(\lambda)$ for chain graphs and $V_{2,3|\emptyset}(\lambda)$ for star graphs by simulation for various values of λ . This means that we sample n points uniformly in the parameter space Ω and count how many of them are $\leq \lambda$. The results for p=6 and p=10 are shown in Figure 6. These are based on a sample size of n=1,000,000. We then fit a linear model

$$\frac{V_{1,p|\emptyset}(\lambda)}{\lambda} \approx C_{p-2}(-\ln \lambda)^{p-2} + C_{p-3}(-\ln \lambda)^{p-3} + \dots + C_0$$

for chain graphs. The curve resulting from the regression estimates is shown in black in Figure 6. The curve resulting from the first-order approximation with the constants computed using Corollary 5.3 is shown in grey in Figure 6. We note that especially for chain graphs, where the true constant in Corollary 5.3 is small, the first order approximation is very bad.

The approximation by regression on the other hand is a fast way to get pretty accurate estimates of all constants. The same was done with star graphs, but with the linear model

$$\frac{V_{2,3|\emptyset}(\lambda)}{\lambda} \approx C_1(-\ln \lambda) + C_0.$$

Figure 6 shows that the first-order approximation is more accurate for stars than for chains.

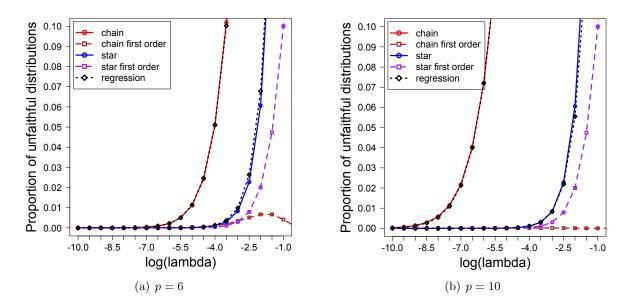


Figure 6: Regression-based asymptotics for chain graphs and stars.

6 Volume inequalities for bias reduction in causal inference

We now discuss the problem of quantifying bias in causal models. Our point of departure is Greenland's paper [7]. In contrast to the previous sections, in the situation discussed here, a large tube volume is in fact desired since it corresponds to small bias. In this section we use the notation $K_{i,j|S}$ for the almost-principal minor $K_{iR,jR}$ of the concentration matrix.

We are interested in estimating the direct effect of an exposure E on a disease outcome D from the partial correlation $\operatorname{corr}(E,D\,|\,S)$, where S is a subset of measurable variables. This partial correlation is a weighted sum over all *open paths* (i.e paths which d-connect E to D) given S (the direct path a_{ED} being just one of them). For estimating the direct effect a_{ED} from $\operatorname{corr}(E,D\,|\,S)$, all open paths other than the direct path are thus considered as bias. We shall analyze two forms of bias which are of particular interest in practice, namely confounding bias and collider-stratification bias. We start by defining collider-stratification bias.

Suppose we are given a DAG G with $D, E \in V$ and there is another node C such that

$$E \to V_1 \to \cdots \to V_s \to C \leftarrow W_1 \leftarrow \cdots \leftarrow W_t \leftarrow D.$$

This says that C is a collider on a path from D to E. Stratifying (i.e. conditioning) on C opens a path between E and D leading to bias when estimating a_{ED} . The partial correlation corresponding to the opened path between E and D is known as *collider-stratification bias*.

Example 6.1. We illustrate collider-stratification bias for the tripartite graph $G = \text{Tripart}_{5,1}$ shown in Figure 7(a). Let node 1 represent the exposure E and node 2 the disease outcome D. In this example, node 5 is a collider C for multiple paths between E and D. When stratifying on C = 5, node E = 1 is d-connected to node D = 2 via the following paths:

$$1 \rightarrow 3 \rightarrow 5 \leftarrow 4 \leftarrow 2, \quad 1 \rightarrow 4 \rightarrow 5 \leftarrow 3 \leftarrow 2, \quad 1 \rightarrow 3 \rightarrow 5 \leftarrow 3 \leftarrow 2, \quad 1 \rightarrow 4 \rightarrow 5 \leftarrow 4 \leftarrow 2.$$

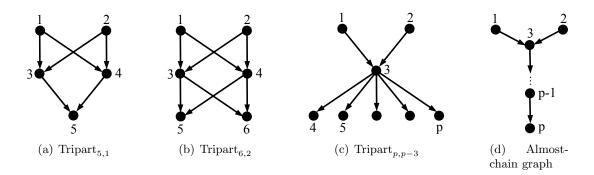


Figure 7: Various tripartite and almost tripartite graphs.

The bias introduced for estimating the direct effect of E on D when conditioning on C is

$$\operatorname{corr}(1, 2 \mid 5) = \frac{a_{13}a_{35}a_{45}a_{24} + a_{14}a_{45}a_{35}a_{23} + a_{13}a_{35}^2a_{23} + a_{14}a_{45}^2a_{24}}{\sqrt{\det(K_{134,134})\det(K_{234,234})}}.$$
 (15)

The numerator $det(K_{134,234})$ is the weighted sum of all open paths between E and D. Similarly, nodes 3 and 4 are colliders for multiple paths. The bias when conditioning on these is

$$\operatorname{corr}(1, 2 \mid 34) = \operatorname{corr}(1, 2 \mid 345) = \frac{a_{13}a_{23} + a_{14}a_{24}}{\sqrt{(a_{13}^2 + a_{14}^2 + 1)(a_{23}^2 + a_{24}^2 + 1)}}.$$
 (16)

Problem 6.2 is about comparing the tube volume for (15) with the tube volume for (16). \Box

A question of practical interest in causal inference is to understand the situations in which stratifying on a collider leads to a particularly large bias. It is widely believed that collider-stratification bias tends to attenuate when it arises from more extended paths (see [2, 7]). What follows is our interpretation of this statement as a precise mathematical conjecture.

Problem 6.2. Let $D, E \in V$ and $C = \{C \in V \mid \exists \text{ path } P \text{ from } E \text{ to } D \text{ with } C \text{ as a collider} \}$. We introduce a partial order on the *collider set* C by setting $C \leq C'$ if all paths on which C is a collider also go through C'. Given subsets $S, S' \subset C$ we set $S \leq S'$ if for all $C \in S$ there exists $C' \in S'$ such that $C \leq C'$. If this holds, then the bias introduced when conditioning on S should be smaller than when conditioning on S'. To make this precise, we conjecture:

$$V_{D,E|S}(\lambda) \ge V_{D,E|S'}(\lambda)$$
 for all $S \le S'$ and all $\lambda \in [0,1]$. (17)

We now study this conjecture for the tripartite graphs $\operatorname{Tripart}_{p,p'}$. Here, it says that the collider-stratification bias introduced when conditioning on the third level $\{p-p'+1,\ldots,p\}$ is in general smaller than when conditioning on the second level of nodes $\{3,\ldots,p-p'\}$, i.e.

$$V_{1,2|p-p'+1,\dots,p}(\lambda) \ge V_{1,2|3,\dots,p-p'}(\lambda).$$
 (18)

This inequality is confirmed by the simulations shown in Figure 8(a). Here p = 5, p' = 2 is shown in red and p = 10, p' = 2 is shown in blue. The solid lines correspond to the volume $V_{1,2|p-p'+1,...,p}(\lambda)$, whereas the dashed lines correspond to the volume $V_{1,2|3,...,p-p'}(\lambda)$.

Going beyond simulations, we now present an algebraic proof of our conjecture when λ is small, for the tripartite graphs in Figure 7(c) where the second level has only one node.

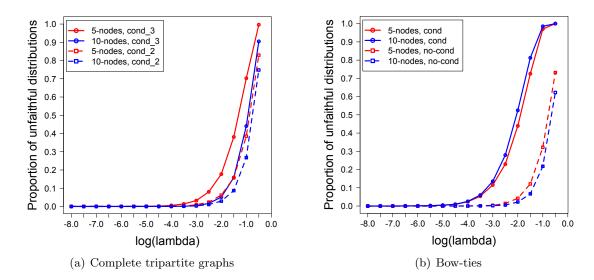


Figure 8: Effect of collider-bias on complete tripartite graphs and bow-ties.

Example 6.3. For $G = \text{Tripart}_{n,n-3}$ the left hand side of (18) is given by

$$\det(K_{1,2|4,5,\dots p}) = a_{13}a_{23}(\sum_{k=4}^{p} a_{3k}^{2}).$$

Depending on the values of p, the corresponding real log canonical threshold is given by

RLCT(1, 2|4,..., p) =
$$\begin{cases} (\frac{1}{2}, 1) & \text{if } p = 4, \\ (1, 3) & \text{if } p = 5, \\ (1, 2) & \text{if } p \ge 6. \end{cases}$$
(19)

For p=4 this was Example 4.7. To prove (19) for $p \geq 5$, we need two ingredients. Firstly, if the polynomial is a product of factors with disjoint variables, then the RLCT is the minimum of the RLCT of the factors, taken with multiplicity (e.g. if the RLCTs are (ℓ, m_1) and (ℓ, m_2) , then the combined RLCT is $(\ell, m_1 + m_2)$, just like in the case of a monomial). Secondly, the RLCT of a sum of squares of ℓ unknowns is equal to ℓ (ℓ , ℓ). We saw this in Example 3.3.

For the right hand side of (18), we condition on node 3. Now, the defining polynomial is

$$\det(K_{1,2|3}) = a_{13}a_{23}.$$

By Proposition 3.5, this has RLCT (1,2), which is larger or equal to all values of (ℓ, m) in (19). To compare the behavior of $V_{1,2|3}(\lambda)$ and $V_{1,2|4,...,p}(\lambda)$ for small λ , we will need to derive the constant C in (7). In Example 8.8, we will show that if $p \geq 6$ and the parameter space is

$$\Omega = \{ a \in \mathbb{R}^{p-1} : |a_{12}| \le 1, |a_{23}| \le 1, a_{34}^2 + \dots + a_{3p}^2 \le 1 \},$$
 (20)

then the asymptotic constants are given by $C_{1,2|3} = 1$ and $C_{1,2|4,...,p} = 2 + 2/(p-5)$. We conclude that $V_{1,2|3}(\lambda) \leq V_{1,2|4,...,p}(\lambda)$ for small values of λ , as conjectured in Problem 6.2. \square

Example 6.4. A slight twist to Example 6.3 is the almost-chain graph shown in Figure 7(d), with edges $E = \{(1,3), (2,3), (3,4), \dots (p-1,p)\}$. For such graphs, Problem 6.2 asks whether

$$V_{1,2|s}(\lambda) \le V_{1,2|t}(\lambda)$$
 if $s \le t$.

This holds for small λ because $\det(K_{1,2|s}) = a_{13}a_{23}\prod_{k=3}^{s-1}a_{k,k+1}^2$. By Proposition 3.5,

RLCT(1,2|s) =
$$\begin{cases} (1,2) & \text{if } s = 3, \\ (\frac{1}{2}, s - 3) & \text{if } s \ge 4, \end{cases}$$

so
$$RLCT(1, 2|s) \ge RLCT(1, 2|t)$$
 for $s \le t$.

In Example 6.3 we resolved Problem 6.2 for tripartite graphs whose middle level consists of one node. We next consider the case $Tripart_{p,1}$ where the third level has one node.

Example 6.5. The graph Tripart_{5,1} shown in Figure 7(a) was discussed in Example 6.1. We focus on the numerators in (15) and in (16). The polynomial (15) will be studied in Example 7.5 where we prove that RLCT(1, 2 | 5) = (1, 3). Using the same method for Tripart_{p,1} gives

$$RLCT(1,2|p) = \begin{cases} (\frac{1}{2},1) & \text{if } p = 4, \\ (1,3) & \text{if } p = 5, \\ (1,2) & \text{if } p \ge 6, \end{cases} \qquad RLCT(1,2|3,\ldots,p-1) = \begin{cases} (1,2) & \text{if } p = 4, \\ (1,1) & \text{if } p \ge 5. \end{cases}$$

Thus, we conclude that $V_{1,2|p}(\lambda) \geq V_{1,2|3,\dots,p-1}(\lambda)$ for small values of $\lambda > 0$.

Example 6.6. For the graph Tripart_{6,2} in Figure 7(b) we check if $V_{1,2|56}(\lambda) \geq V_{1,2|34}(\lambda)$ for small λ . As before, RLCT(1,2|3,4) = (1,1), but a hard computation using the tools of Section 7 reveals that now RLCT(1,2|5,6) = (1,1). Thus, knowledge of the RLCT is not sufficient to establish (17). What is needed is a finer analysis along the lines of Section 8.

The second form of bias studied by Greenland [7] is confounder bias. In the context of a directed graphical model G, a confounder for the effect of E on D is a node C such that

$$E \leftarrow V_1 \leftarrow \cdots \leftarrow V_s \leftarrow C \rightarrow W_1 \rightarrow \cdots \rightarrow W_t \rightarrow D.$$

The partial correlation introduced by the path from E to D passing through C is referred to as *confounder bias*. In such situations, stratifying on C blocks the path between E and D (i.e. C d-separates E from D) and therefore corresponds to bias removal.

In certain graphs, such as the bow-tie example in [7], there are variables where stratifying removes confounder bias but at the same time introduces collider-stratification bias. For instance, consider the graph $G = \text{Bow}_5$, where node 4 corresponds to exposure E and node 5 corresponds to disease outcome D. Then conditioning on node 3 blocks the paths

$$4 \leftarrow 3 \rightarrow 5, \quad 4 \leftarrow 1 \rightarrow 3 \rightarrow 5, \quad 4 \leftarrow 3 \leftarrow 2 \rightarrow 5,$$
 (21)

and therefore reduces confounder bias, but opens the path

$$4 \leftarrow 1 \rightarrow 3 \leftarrow 2 \rightarrow 5 \tag{22}$$

and therefore introduces collider-stratification bias. This trade-off is of particular interest in situations where one cannot condition on 1 and 2, for example because these variables were unmeasured. It is believed that in such examples the bias removed by conditioning on the confounders is larger than the collider-stratification bias introduced and one should therefore stratify. We translate this statement into the following mathematical problem:

Problem 6.7. Let $D, E \in V$ and we denote by \mathcal{D} the confounder-collider subset, i.e.

$$\mathcal{D} = \mathcal{C} \cap \{C \in V \mid \exists \text{ path } \pi \text{ from } E \text{ to } D \text{ having } C \text{ as a confounder}\}.$$

We conjecture the following inequality for the relevant tube volumes:

$$V_{D,E|S}(\lambda) \geq V_{D,E|\emptyset}(\lambda),$$
 for all $S \subset \mathcal{D}$ and all $\lambda \in [0,1]$.

This conjectural inequality is interesting for the bow-tie graphs Bow_p . It means that conditioning on the nodes in the second level reduces bias since the bias removed by conditioning on the confounders is larger than the collider-stratification bias introduced by conditioning:

$$V_{p-1,p|3,...,p-2}(\lambda) \ge V_{p-1,p|\emptyset}(\lambda).$$
 (23)

This is confirmed by our simulations in Figure 8(b), for p=5 in red and p=10 in blue. The solid line corresponds to the volume $V_{p-1,p|3,...,p-2}(\lambda)$ and the dashed line corresponds to $V_{p,p-1|\emptyset}(\lambda)$. In the following example we prove the inequality (23) for p=5 and small $\lambda > 0$.

Example 6.8. Let $G = \text{Bow}_5$ as in Figure 1(d). The left hand side (23) is represented by

$$\det(K_{4,5|3}) = a_{13}a_{14}a_{23}a_{25}.$$

This monomial is the path in (22). The corresponding real log canonical threshold is (1,4). The polynomial representing the right hand side (23) is a weighted sum of the paths in (21):

$$\det(K_{4,5|\emptyset}) = a_{34}a_{35}(1 + a_{13}^2 + a_{23}^2) + a_{23}a_{25}a_{34} + a_{13}a_{14}a_{35}.$$

We derive its real log canonical threshold using the blowups described in Section 7. We find that it is (1,1). Since (1,4)<(1,1), we conclude $V_{4,5|3}(\lambda) \geq V_{4,5|\emptyset}(\lambda)$.

7 Normal crossing and blowing up

In this section we develop more refined techniques for computing real log canonical thresholds. The following theorem combines the monomial case of Proposition 3.5 with the smooth case of Proposition 3.6. As promised in Section 3, this furnishes the proofs for these two propositions.

Theorem 7.1. Suppose $\varphi(\omega) = \omega_1^{\tau_1} \cdots \omega_d^{\tau_d}$ and $f(\omega) = \omega_1^{\kappa_1} \cdots \omega_r^{\kappa_r} g(\omega)$ where $\kappa_1, \ldots, \kappa_r$ are non-zero and the hypersurface $g(\omega) = 0$ is smooth and normal crossing (see definition below) with $\omega_1, \ldots, \omega_r$. Let ω_0 denote the function g and let $\kappa_0 = 1, \tau_0 = 0$. Define

$$\ell = \min_{i \in \mathcal{I}} \frac{\tau_i + 1}{\kappa_i}, \quad \mathcal{J} = \underset{i \in \mathcal{I}}{\operatorname{argmin}} \frac{\tau_i + 1}{\kappa_i}, \quad m = |\mathcal{J}|,$$

where \mathcal{I} is the set of all indices $0 \leq i \leq r$ such that ω_i has a zero in Ω . Then we have

$$RLCT_{\Omega}(f;\varphi) = (\ell, m),$$

provided the equations $\omega_i = 0$ for $i \in \mathcal{J}$ have a solution in the interior of Ω .

The normal crossing hypothesis in Theorem 7.1 means that the system

$$f = \omega_1 \frac{\partial f}{\partial \omega_1} = \dots = \omega_r \frac{\partial f}{\partial \omega_r} = \frac{\partial f}{\partial \omega_{r+1}} = \dots = \frac{\partial f}{\partial \omega_d} = 0$$

does not have any solutions in Ω . See [12] to learn more about normal crossing singularities. We begin with a technical lemma establishing that the RLCT can be computed locally.

Lemma 7.2. For every $x \in \Omega$, there exists a neighborhood $\Omega_x \subset \Omega$ of x such that

$$RLCT_{\Omega_x}(f;\varphi) = RLCT_U(f;\varphi)$$

for all neighborhoods $U \subset \Omega_x$ of x. Moreover,

$$RLCT_{\Omega}(f;\varphi) = \min_{x} RLCT_{\Omega_{x}}(f;\varphi)$$

where we take the minimum over all x in the real analytic hypersurface $\{\omega \in \Omega : f(\omega) = 0\}$.

Proof. This comes from [14, Lemma 3.8, Proposition 3.9].

Proof of Theorem 7.1. Lemma 7.2 states that $\mathrm{RLCT}_{\Omega}(f;\varphi)$ is the minimum of $\mathrm{RLCT}_{\Omega_x}(f;\varphi)$ as x varies over Ω . Writing each subset Ω_x as $R_x \cap \Omega$ where R_x is a sufficiently small neighborhood of x in \mathbb{R}^d , we claim that $\mathrm{RLCT}_{\Omega}(f;\varphi) = \min_{x \in \Omega} \mathrm{RLCT}_{R_x}(f;\varphi)$ if this minimum is attained in the interior of Ω . Indeed, for x in the interior of Ω , we get $\mathrm{RLCT}_{R_x}(f;\varphi) = \mathrm{RLCT}_{\Omega_x}(f;\varphi)$. Otherwise, the volume of $\{\omega \in \Omega_x : f(\omega) \leq \lambda\}$ is less than that of $\{\omega \in R_x : f(\omega) \leq \lambda\}$ for all λ . Hence $\mathrm{RLCT}_{R_x}(f;\varphi) \leq \mathrm{RLCT}_{\Omega_x}(f;\varphi)$, and the claim follows.

Now to prove Theorem 7.1, it suffices to show that for each $x \in \Omega$ we have

$$RLCT_{R_x}(f;\varphi) = \left(\min_{i \in \mathcal{I}_x} \frac{\tau_i + 1}{\kappa_i}, \left| \underset{i \in \mathcal{I}_x}{\operatorname{argmin}} \frac{\tau_i + 1}{\kappa_i} \right| \right)$$

where \mathcal{I}_x is the set of all indices $0 \leq i \leq r$ that satisfies $\omega_i(x) = 0$. Without loss of generality, suppose $x = (x_1, \ldots, x_d)$ where $x_1 = \cdots = x_s = 0$ and x_{s+1}, \ldots, x_r are nonzero. If $g(x) \neq 0$, we may divide $f(\omega)$ by $g(\omega)$ without changing the RLCT in a sufficiently small neighborhood R_x of x. The RLCT of the remaining monomial is determined by [14, Proposition 3.7]. Now, let us suppose g(x) = 0. Because $g(\omega)$ is normal crossing with $\omega_1, \ldots, \omega_r$, one of the derivatives $\partial g/\partial \omega_j$ must be nonzero at x for some $s+1 \leq j \leq d$. We assume R_x is sufficiently small so that this derivative and $\omega_{s+1}, \ldots, \omega_r$ do not vanish. Consider the map $\sigma: R_x \to \mathbb{R}^d$ given by

$$\sigma_i(\omega) = g(\omega), \quad \sigma_i(\omega) = \omega_i \quad \text{for } i \neq j.$$

The Jacobian matrix of σ is nonsingular, so this map is an isomorphism onto its image. Set $U = \mu(R_x)$ and $\rho = \sigma^{-1} : U \to R_x$. Then, for all $\mu \in U$, we have

$$(f \circ \rho)(\mu) = \mu_1^{\kappa_1} \cdots \mu_s^{\kappa_s} \mu_j \cdot a(\mu)$$
 and $(\varphi \circ \rho)(\mu) = \mu_1^{\tau_1} \cdots \mu_s^{\tau_s} \cdot b(\mu),$

where the factors $a(\mu)$ and $b(\mu)$ do not vanish in U. By using the chain rule [14, Proposition 4.6], we get $\mathrm{RLCT}_{R_x}(f;\varphi) = \mathrm{RLCT}_U(f \circ \rho; \varphi \circ \rho)$. The latter RLCT can be computed once again by dividing out the nonvanishing factors and applying [14, Proposition 3.7].

The hypersurface $\{f(\omega) = 0\}$ may not satisfy the hypothesis in Theorem 7.1. In that case, we can try to simplify its singularities via a change of variables $\rho: U \to \Omega$. With some luck, the transformed hypersurface $\{(f \circ \rho)(\mu) = 0\}$ will be described locally by monomials and the RLCT can be computed using Theorem 7.1. More precisely, let U be a d-dimensional real analytic manifold and $\rho: U \to \Omega$ a real analytic map that is proper, i.e. the preimage of any compact set is compact. Then ρ desingularizes $f(\omega)$ if it satisfies the following conditions:

- 1. The map ρ is an isomorphism outside the variety $\{\omega \in \Omega : f(\omega) = 0\}$.
- 2. Given any $y \in U$, there exists a local chart with coordinates μ_1, \ldots, μ_d such that

$$(f \circ \rho)(\mu) = \mu_1^{\kappa_1} \cdots \mu_d^{\kappa_d} \cdot a(\mu), \quad \det \partial \rho(\mu) = \mu_1^{\tau_1} \cdots \mu_d^{\tau_d} \cdot b(\mu)$$

where det $\partial \rho$ is the Jacobian determinant, the exponents κ_i , τ_i are nonnegative integers and the real analytic functions $a(\mu)$, $b(\mu)$ do not vanish at y.

If such a desingularization exists, then we may apply ρ to the volume function (7) to calculate the RLCT. Care must be taken to multiply the measure φ with the Jacobian determinant $|\det \partial \rho|$ in accordance with the change-of-variables formula for integrals.

Hironaka's celebrated theorem on the resolution of singularities [9, 10] guarantees that such a desingularization exists for all real analytic functions $f(\omega)$. The proof employs transformations known as blowups to simplify the singularities. We now describe the blowup $\rho: U \to \mathbb{R}^d$ of the origin in \mathbb{R}^d . The manifold U can be covered by local charts U_1, \ldots, U_d such that each chart is isomorphic to \mathbb{R}^d and each restriction $\rho_i: U_i \to \mathbb{R}^d$ is the monomial map

$$(\mu_1, \dots, \mu_{i-1}, \xi, \mu_{i+1}, \dots, \mu_d) \mapsto (\xi \mu_1, \dots, \xi \mu_{i-1}, \xi, \xi \mu_{i+1}, \dots, \xi \mu_d).$$

Here, the coordinate hypersurface $\xi = 0$, also called the *exceptional divisor*, runs through all the charts. If the origin is locally the intersection of many smooth hypersurfaces with distinct tangent hyperplanes, then these hypersurfaces can be separated by blowing up the origin [9].

Example 7.3. Consider the curve $\{f(x,y) = xy(x+y)(x-y) = 0\}$ in Figure 3(d). To resolve this singularity, we blow up the origin. In the first chart, the map is $\rho_1 : (\xi, y_1) \mapsto (\xi, \xi y_1)$, so

$$f \circ \rho_1 = \xi^4 y_1 (1 + y_1) (1 - y_1)$$
 and $\det \partial \rho_1 = \xi$.

The lines $\{y=0\}$, $\{x+y=0\}$ and $\{x-y=0\}$ are transformed to $\{y_1=0\}$, $\{y_1=-1\}$ and $\{y_1=1\}$ respectively in this chart, thereby separating them. The line $\{x=0\}$ does not show up here, but it appears as $\{x_1=0\}$ in the second chart, where $\rho_2: (x_1,\xi) \mapsto (\xi x_1,\xi)$ and

$$f \circ \rho_2 = \xi^4 x_1(x_1 + 1)(x_1 - 1), \quad \det \partial \rho_2 = \xi.$$

Since the curve $\{(1+y_1)(1-y_1)=0\}$ is normal crossing with ξ, y_1 in the first chart, we can now apply Theorem 7.1. The chain rule [14, Proposition 4.6] shows that $RLCT_{\Omega}(f;1)$ is the minimum of $RLCT_{U_i}(f \circ \rho_i; \det \partial \rho_i)$ for i=1,2. In both charts, this RLCT equals $(\frac{1}{2},1)$. \square

Example 7.4. Let p = 4 and G be the almost complete DAG with $a_{13} = 0$. We consider the conditional independence statement $1 \perp \!\!\! \perp 3 \mid 4$. The correlation hypersurface is defined by

$$f = \det(K_{12,23}) = a_{14}a_{23}^2a_{34} + a_{14}a_{23}a_{24} + a_{12}a_{24}a_{34} - a_{12}a_{23} + a_{14}a_{34}.$$

The real singular locus is a line in the parameter space \mathbb{R}^5 , since $\operatorname{Singu}_{1,3,2} = \langle a_{12}, a_{14}, a_{23}, a_{34} \rangle$. Blowing up this line in \mathbb{R}^5 creates four charts U_1, U_2, U_3, U_4 . For instance, the first chart has

$$\rho_1: U_1 \to \mathbb{R}^5, \ (\xi, \mu_{14}, \mu_{23}, a_{24}, \mu_{34}) \mapsto (\xi, \xi \mu_{14}, \xi \mu_{23}, a_{24}, \xi \mu_{34}), \ \det \partial \rho_1 = \xi^3.$$

Then f transforms to $f \circ \rho_1 = \xi^2 \cdot g$ where $g = \mu_{14}\mu_{23}^2\mu_{34}\xi^2 + \mu_{14}\mu_{34} + \mu_{14}\mu_{23}a_{24} + \mu_{34}a_{24} - \mu_{23}$. The hypersurface $\{g = 0\}$ has no real singularities, so it is smooth in U_1 . We can thus apply Theorem 7.1 with $\mathcal{I} = \{0,1\}$ to find $\mathrm{RLCT}_{U_1}(\xi^2 \cdot g, \xi^3) = (1,1)$. The behavior is the same on U_2, U_3 and U_4 , and we conclude that $\mathrm{RLCT}(1,3|4) = (\ell,m) = (1,1)$. This example is one of the 12 cases that were labeled as "Blowup" in Table 2. The other 11 cases are similar. \square

Example 7.5. In Example 6.5 we claimed that RCLT(1, 2 | 5) = (1, 3) for $G = Tripart_{5,1}$. We now prove this claim by using the blowing up method. The polynomial in question is

$$f = \det(K_{1,2|5}) = (a_{13}a_{35} + a_{14}a_{45})(a_{23}a_{35} + a_{24}a_{45}).$$

The singular locus of the hypersurface $\{f=0\}$ is given by

$$\operatorname{Singu}_{1,2,34} = \langle a_{35}, a_{45} \rangle \cap \langle 2 \times 2 \text{-minors of} \begin{pmatrix} a_{13} & a_{23} & a_{45} \\ a_{14} & a_{24} & -a_{35} \end{pmatrix} \rangle.$$

We blow up the linear subspace $\{a_{35}=a_{45}=0\}$ in \mathbb{R}^6 . This creates two charts. The map for the first chart is $\rho_1:(a_{13},a_{14},a_{23},a_{24},\xi,\mu_{45})\mapsto(a_{13},a_{14},a_{23},a_{24},\xi,\xi\mu_{45})$. This map gives

$$f \circ \rho_1 = \xi^2 (a_{13} + a_{14}\mu_{45})(a_{23} + a_{24}\mu_{45}), \quad \det \partial \rho_1 = \xi.$$

Now, by setting $a_{13} = x - a_{14}\mu_{45}$ and $a_{23} = y - a_{24}\mu_{45}$, the transformed function $f \circ \rho_1$ is the monomial $\xi^2 xy$. Then Theorem 7.1 can be employed to evaluate $\mathrm{RLCT}_{U_1}(\xi^2 xy, \xi) = (1, 3)$. The calculation in the second chart is completely analogous.

The same approach as in Example 7.5 can be applied to the polynomial $f = \det(K_{1,2|56})$ in Example 2.1. A lengthy calculation, involving many charts and multiple blowups, eventually reveals that $G = \text{Tripart}_{6,2}$ satisfies RCLT(1,2|56) = (1,1). This was stated in Example 6.6.

8 Computing the constants

We now describe a method for finding the constant C in the formula $V(\lambda) \approx C\lambda^{-\ell}(-\ln \lambda)^{m-1}$ in (7). The two theorems in this section are new. They extend the results of Greenblatt [6] and Lasserre [13] on the volumes of sublevel sets. We begin by showing that the constant C is a function of the highest order term in the Laurent expansion of the zeta function of f.

Lemma 8.1. Given real analytic functions $f, \varphi : \Omega \to \mathbb{R}$, consider the Laurent expansion of

$$\zeta(z) := \int_{\Omega} |f(\omega)|^{-z} \varphi(\omega) d\omega = \frac{a_{\ell,m}}{(\ell - z)^m} + \frac{a_{\ell,m-1}}{(\ell - z)^{m-1}} + \cdots$$

where ℓ is the smallest pole and m its multiplicity. Then, asymptotically as λ tends to zero,

$$V(\lambda) := \int_{|f(\omega)| < \lambda} \varphi(\omega) d\omega \approx \frac{a_{\ell,m}}{\ell(m-1)!} \lambda^{\ell} (-\ln \lambda)^{m-1}.$$

Proof. According to the proof of [18, Theorem 7.1], the volume function $V(\lambda)$ equals $\int_0^{\lambda} v(s)ds$ where $v(s) = \int_{\Omega} \delta(s - f(\omega)) \varphi(\omega) d\omega$ is the state density function and δ is the delta function. Now, using the proof of [14, Theorem 3.16], we obtain

$$v(s) = \frac{a_{\ell,m}}{(m-1)!} s^{\ell-1} (-\ln s)^{m-1} + o(s^{\ell-1} (-\ln s)^{m-1}) \quad \text{as } s \to 0.$$

Here we used the little-o notation. Finally, using integration by parts, we find that

$$V(\lambda) = \frac{a_{\ell,m}}{\ell(m-1)!} \lambda^{\ell} (-\ln \lambda)^{m-1} + o(\lambda^{\ell} (-\ln \lambda)^{m-1}) \quad \text{as } \lambda \to 0.$$

Example 8.2. In Example 3.3, we saw that the volume of the d-dimensional ball defined by $|\omega_1^2 + \cdots + \omega_d^2| \leq \lambda$ is equal to $V(\lambda) = C\lambda^{-d/2}$ for some positive constant C. We here show how to compute that constant using asymptotic methods. By Lemma 8.1, $C = 2\alpha/(d 2^d)$ where α is the coefficient of $(d/2 - z)^m$ in the Laurent expansion of the zeta function

$$\zeta(z) = \int_{\mathbb{R}^d} |\omega_1^2 + \ldots + \omega_d^2|^{-z} d\omega.$$

Computing this Laurent coefficient from first principles is not easy. Instead, we derive α using the asymptotic theory of Laplace integrals. The connection between such integrals and volume functions was alluded to in Definition 3.2. By [14, Proposition 5.2], the Laplace integral

$$Z(N) = \int_{\mathbb{R}^d} e^{-N(\omega_1^2 + \dots + \omega_d^2)} d\omega$$

is asymptotically $\alpha \Gamma(\frac{d}{2})N^{-d/2}$ for large N. But this Laplace integral also decomposes as

$$Z(N) = \int_{\mathbb{R}} e^{-N\omega_1^2} d\omega_1 \cdots \int_{\mathbb{R}} e^{-N\omega_d^2} d\omega_d = (\sqrt{\pi}N^{-1/2})^d,$$

where each factor is the classical Gaussian integral. Solving for α leads to the formula

$$C = \frac{\pi^{d/2}}{2^d \cdot \Gamma(\frac{d}{2}) \cdot \frac{d}{2}} = \frac{\pi^{d/2}}{2^d \cdot \Gamma(\frac{d}{2} + 1)}.$$

In Section 3, we saw how the RLCTs of smooth hypersurfaces and of hypersurfaces defined by monomial functions can be computed. The following two theorems and their accompanying examples demonstrate how the asymptotic constant C can also be evaluated in those instances.

Theorem 8.3. Let $\{f = 0\}$ be a smooth hypersurface and let $\varphi : \Omega \to \mathbb{R}$ be positive. Suppose $\partial f/\partial \omega_1$ is nonvanishing in Ω . Let W be the projection of the hypersurface $\{f = 0\} \subset \Omega$ onto the subspace $\{(\omega_2, \ldots, \omega_d) \in \mathbb{R}^{d-1}\}$ and let $\rho : \Omega \to \mathbb{R}^d$ be the map $\omega \mapsto (f(\omega), \omega_2, \ldots, \omega_d)$. If the boundary of Ω intersects transversally with the hypersurface $\{f = 0\}$, then

$$V(\lambda) := \int_{\{\omega \in \Omega : |f(\omega)| \le \lambda\}} \varphi(\omega) d\omega \approx C\lambda$$

asymptotically (as $\lambda \to 0$), where

$$C = 2 \int_{W} \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(0, \omega_2, \dots, \omega_d) \, d\omega_2 \cdots d\omega_d.$$

Proof. The asymptotics of the volume $V(\lambda)$ depends only on the region $\{\omega \in \Omega : |f(\omega)| \leq \lambda\}$. So we may assume that Ω is a small neighborhood of the hypersurface $\{f(\omega) = 0\}$. As we saw in the proof of Theorem 7.1, the map ρ is an isomorphism onto its image. Thus after changing variables, the zeta function associated to $V(\lambda)$ becomes

$$\zeta(z) = \int_{\rho(\Omega)} |f|^{-z} \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(f, \omega_2, \dots, \omega_d) \, df \, d\omega_2 \cdots d\omega_d
= \int_W \int_{\varepsilon_1(\omega_2, \dots, \omega_d)}^{\varepsilon_2(\omega_2, \dots, \omega_d)} |f|^{-z} \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(f, \omega_2, \dots, \omega_d) \, df \, d\omega_2 \cdots d\omega_d.$$

Here, the lower and upper limits $\varepsilon_1, \varepsilon_2$ are nonzero because the boundary of Ω is transversal to the hypersurface. By substituting the Taylor series

$$\frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(f, \omega_2, \dots, \omega_d) = \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(0, \omega_2, \dots, \omega_d) + O(f)$$

and the exponential series $\varepsilon_2^{1-z}=1+O(1-z),$ we get the Laurent expansion

$$\int_0^{\varepsilon_2} |f|^{-z} \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(f, \omega_2, \dots, \omega_d) df = \left[\frac{|f|^{1-z}}{1-z} \cdot \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(0, \omega_2, \dots, \omega_d) \right]_0^{\varepsilon_2} + \dots$$

$$= \frac{1}{1-z} \cdot \frac{\varphi}{|\partial_{\omega_1} f|} \circ \rho^{-1}(0, \omega_2, \dots, \omega_d) + \dots$$

The same is true for the integral from ε_1 to 0. The result now follows from Lemma 8.1.

Example 8.4. By Theorem 4.1, all conditional independence statements in small complete graphs lead to smooth hypersurfaces. Here we analyze the statement $1 \perp \!\!\! \perp 2 \mid 3$ in the complete 3-node DAG. This example was studied in [17, §2]. The corresponding partial correlation is

$$corr(1, 2 | 3) = -\frac{a_{13}a_{23} - a_{12}}{\sqrt{1 + a_{23}^2}\sqrt{1 + a_{12}^2 + a_{13}^2}}.$$

This partial correlation hypersurface lives in \mathbb{R}^3 and it is depicted in [17, Figure 2(b)].

We apply Theorem 8.3 by setting $\Omega := [-1,1]^3$, $f := \operatorname{corr}(1,2|3)$ and $\varphi := 1/2^3$, the uniform distribution on Ω . We choose ω_1 to be a_{12} . Then $\rho^{-1}(0,a_{13},a_{23}) = (a_{13}a_{23},a_{13},a_{23})$.

The projection W of the surface $\{a_{12} = a_{13}a_{23}\}$ onto $\{(a_{13}, a_{23}) \in [-1, 1]^2\}$ is the whole square $[-1, 1]^2$. The formula for the constant C now simplifies to

$$C = \frac{1}{4} \int_{-1}^{1} \int_{-1}^{1} \sqrt{1 + a_{13}^2} \sqrt{1 + a_{13}^2 + a_{13}^2 a_{23}^2} \ da_{13} \ da_{23} \approx 5.4829790759.$$

This two-dimensional integral was evaluated numerically using Mathematica.

We now come to the monomial case that was discussed in Theorem 7.1.

Theorem 8.5. Let $f(\omega) = \omega_1^{\kappa_1} \cdots \omega_d^{\kappa_d} g(\omega)$ where $g: \Omega \to \mathbb{R}$ has no real zeros. Let $\varphi: \Omega \to \mathbb{R}$ be positive. Suppose that $1/\ell = \kappa_1 = \cdots = \kappa_m > \kappa_{m+1} \geq \cdots \geq \kappa_d$ and that the boundary of Ω is transversal to the subspace L defined by $\omega_1 = \cdots = \omega_m = 0$. Let $\bar{\omega}$ and $\bar{\kappa}$ denote the vectors $(\omega_{m+1}, \ldots, \omega_d)$ and $(\kappa_{m+1}, \ldots, \kappa_d)$ respectively. Then

$$V(\lambda) := \int_{\{\omega \in \Omega: |f(\omega)| \le \lambda\}} \varphi(\omega) d\omega \approx C\lambda^{\ell} (-\ln \lambda)^{m-1}$$

asymptotically as λ tends to zero where

$$C = \frac{(2\ell)^m}{\ell(m-1)!} \int_{\Omega \cap L} \bar{\omega}^{-\ell\bar{\kappa}} g(0,\dots,0,\bar{\omega})^{-\ell} \varphi(0,\dots,0,\bar{\omega}) d\bar{\omega}. \tag{24}$$

Proof. Let us suppose for now that Ω is the hypercube $[0,\varepsilon]^d$. Because $f(\omega)$ is nondegenerate with respect to its Newton polyhedron [14, §4.2.1], we may apply the last formula in the proof of [14, Theorem 5.11], with $\rho_{\mathcal{F}}$ equal to the identity map. From this we obtain the desired Laurent coefficient $a_{\ell,m}$ of the zeta function, namely

$$a_{\ell,m} = \prod_{i=1}^{m} \frac{1}{\kappa_i} \cdot \int_{[0,\varepsilon]^{d-m}} \bar{\omega}^{-\ell\bar{\kappa}} g(0,\ldots,0,\bar{\omega})^{-\ell} \varphi(0,\ldots,0,\bar{\omega}) d\bar{\omega}.$$

More generally, when the boundary of Ω is transversal to the subspace L, we decompose Ω into small neighborhoods which are isomorphic to orthants. Summing up over these orthants and applying Lemma 8.1 gives the desired result.

Remark 8.6. We revisit the planar tubes shown in Figures 3(a)-3(c). Using the formula (24) in Theorem 8.5, one can easily check the constants C we saw in Example 3.1, namely

$$C = \begin{cases} 1 & \text{for } f(x,y) = x, \\ 1 & \text{for } f(x,y) = xy, \\ 3 & \text{for } f(x,y) = x^2 y^3. \end{cases}$$

Example 8.7. We apply Theorem 8.5 to find the constants in Corollary 5.3 for chains and stars. In both cases we set $\Omega = [-1,1]^{p-1}$ and $\varphi = 2^{1-p}$. For chains we have $(\ell,m) = (1,p-1)$ and L is the subspace $a_{12} = \cdots = a_{p-1,p} = 0$. Then the integral in (24) is the evaluation of the denominator of (13) at the origin multiplied by φ , so $C_{\text{chain}} = 1/(p-2)!$ as claimed.

For stars we have $(\ell, m) = (1, 2)$ and one can check that $\operatorname{corr}(i, j) \geq \operatorname{corr}(i, j|S)$ for all triples i, j, S. Hence, $V_G(\lambda)$ is the volume of the union of the tubes $\{|\operatorname{corr}(i, j)| \leq \lambda\}$ where 1 < i < j. By applying formula (24) to the function (14), the asymptotic volume of each tube computes to $\lambda(-\ln \lambda)$. Meanwhile, as $\lambda \to 0$, the volumes of the intersections of these tubes become negligible. Therefore C_{star} equals $\binom{p-1}{2}$, the number of pairs 1 < i < j.

Example 8.8. We compute the constant C of the volume $V_{1,2|4,...,p}(\lambda)$ for $G = \text{Tripart}_{p,p-3}$ as in Example 6.3. Let $p \geq 6$ and Ω the parameter space (20). By symmetry, it suffices to study the intersection Ω_+ of Ω with the positive orthant. Let $g(\bar{a}) = \sum_{k=4}^p a_{3k}^2$. To apply Lemma 8.1, we need to derive the coefficient of $(1-z)^2$ in the Laurent expansion of

$$\zeta(z) = \int_{\Omega_{\perp}} \left(a_{13} a_{23} \, \frac{g(\bar{a})}{h(a)} \right)^{-z} da, \quad \text{where } h(a) = \sqrt{1 + g(\bar{a})(a_{13}^2 + 1)} \cdot \sqrt{1 + g(\bar{a})(a_{23}^2 + 1)}.$$

The Taylor series expansion of the integrand about $a_{13} = a_{23} = 0$ is

$$\left(a_{13}a_{23}\frac{g(\bar{a})}{h(a)}\right)^{-z} = (a_{13}a_{23})^{-z} \left[\left(\frac{g(\bar{a})}{1+g(\bar{a})}\right)^{-z} + O(a_{13}) + O(a_{23}) \right].$$

The higher order terms in the Taylor expansion contribute larger poles to the zeta function $\zeta(z)$, so we only need to study the Laurent series

$$\int_{\Omega_{+}} (a_{13}a_{23})^{-z} \left(\frac{g(\bar{a})}{1+g(\bar{a})}\right)^{-z} da_{13}da_{23}d\bar{a} = \frac{1}{(1-z)^{2}} \int_{\{g(\bar{a})\leq 1\}} \left(\frac{g(\bar{a})}{1+g(\bar{a})}\right)^{-1} d\bar{a} + \cdots$$

By using the Lebesgue probability measure on the ball $\{g(\bar{a}) \leq 1\}$ and substituting spherical coordinates for the integration, we find that the above coefficient of $(1-z)^2$ equals

$$1 + \int_{\{q(\bar{a}) \le 1\}} \frac{1}{g(\bar{a})} d\bar{a} = 2 + \frac{2}{p-5}.$$

This is the constant $C_{1,2|4,\ldots,p}$ needed for the bias reduction analysis in Example 6.3.

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